



Full wwPDB NMR Structure Validation Report ⓘ

Mar 25, 2026 – 04:54 PM UTC

PDB ID : 2KI4 / pdb_00002ki4
Title : FGF1-S100A13 complex structure: key component in non-classical path way of FGF1
Authors : Krishna, S.M.; Rani, S.G.; Yu, C.
Deposited on : 2009-04-27

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

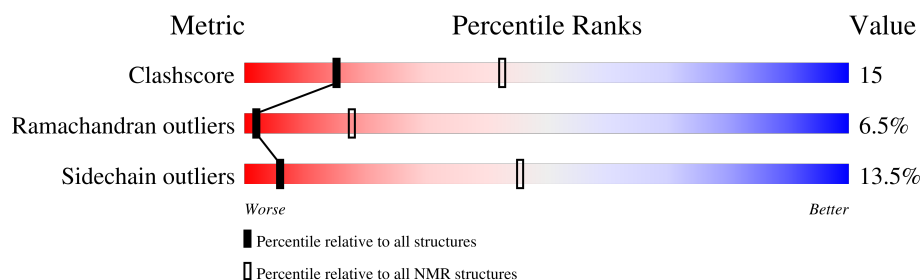
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	133	51% 40% 8% .
1	D	133	53% 38% 8% .
2	B	98	69% 19% . 9%
3	C	98	60% 27% . 12%

2 Ensemble composition and analysis

This entry contains 20 models. Model 7 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:1-A:133, B:8-B:96, C:12-C:93, C:95-C:98, D:1-D:133 (441)	0.27	7

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 2 clusters and 1 single-model cluster was found.

Cluster number	Models
1	1, 2, 3, 6, 7, 8, 9, 10, 11, 12, 13, 14, 16, 17, 18, 19, 20
2	5, 15
Single-model clusters	4

3 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7505 atoms, of which 3764 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Heparin-binding growth factor 1.

Mol	Chain	Residues	Atoms						Trace
1	A	133	Total	C	H	N	O	S	0
			2113	671	1049	185	204	4	
1	D	133	Total	C	H	N	O	S	0
			2113	671	1049	185	204	4	

- Molecule 2 is a protein called Protein S100-A13.

Mol	Chain	Residues	Atoms						Trace
2	B	98	Total	C	H	N	O	S	0
			1640	512	834	136	156	2	

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	97	DLY	LYS	SEE REMARK 999	UNP Q99584
B	98	DLY	LYS	SEE REMARK 999	UNP Q99584

- Molecule 3 is a protein called Protein S100-A13.

Mol	Chain	Residues	Atoms						Trace
3	C	98	Total	C	H	N	O	S	0
			1639	512	832	136	157	2	

There is a discrepancy between the modelled and reference sequences:

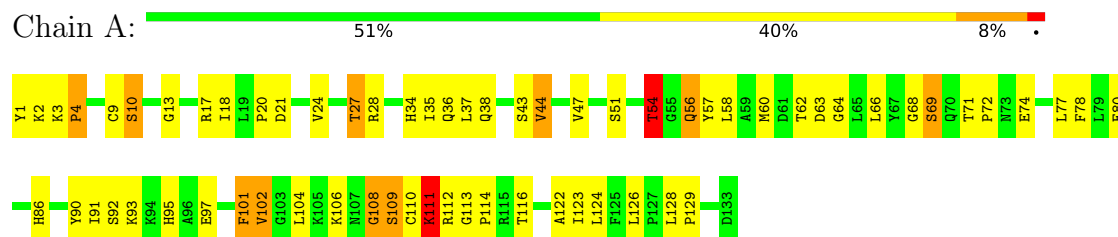
Chain	Residue	Modelled	Actual	Comment	Reference
C	94	DLY	LYS	SEE REMARK 999	UNP Q99584

4 Residue-property plots [i](#)

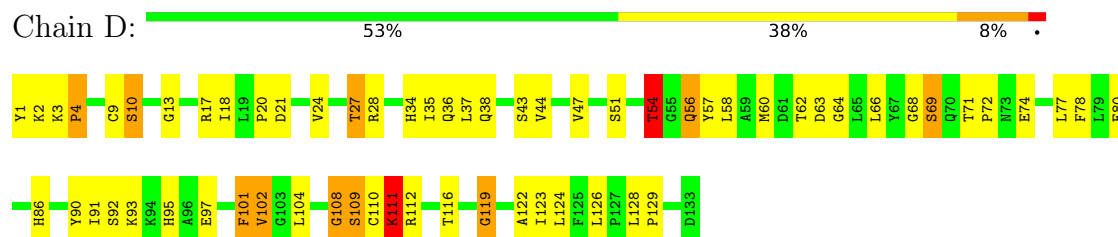
4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

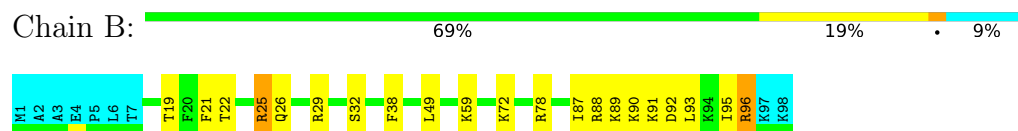
- Molecule 1: Heparin-binding growth factor 1



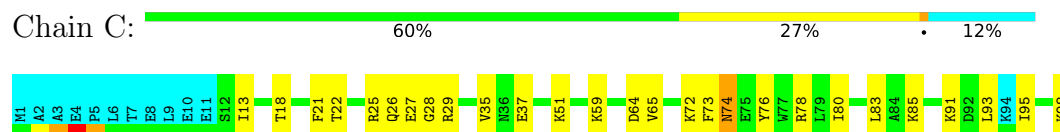
- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13



- Molecule 3: Protein S100-A13

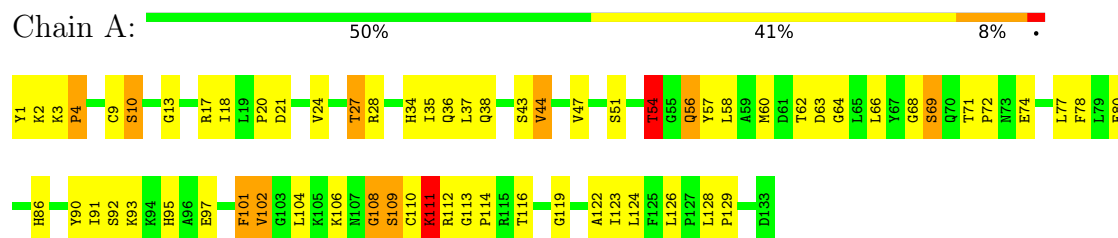


4.2 Scores per residue for each member of the ensemble

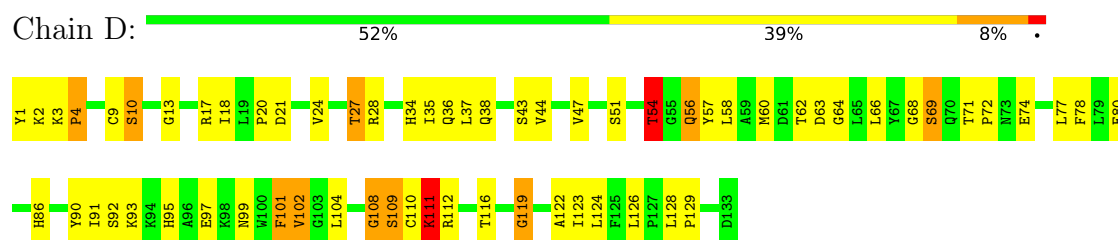
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

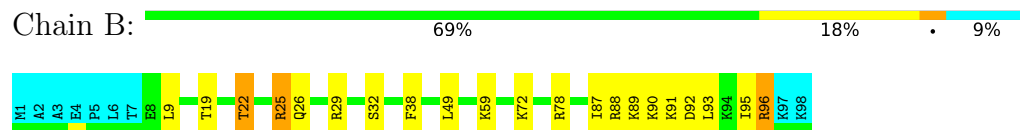
- Molecule 1: Heparin-binding growth factor 1



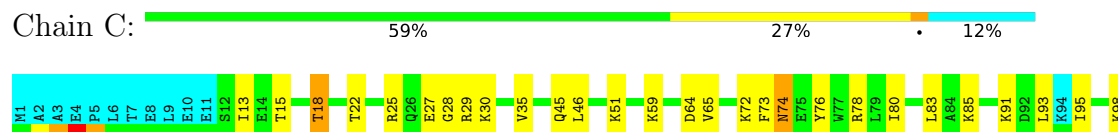
- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13

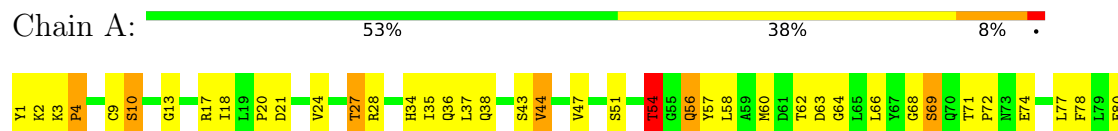


- Molecule 3: Protein S100-A13



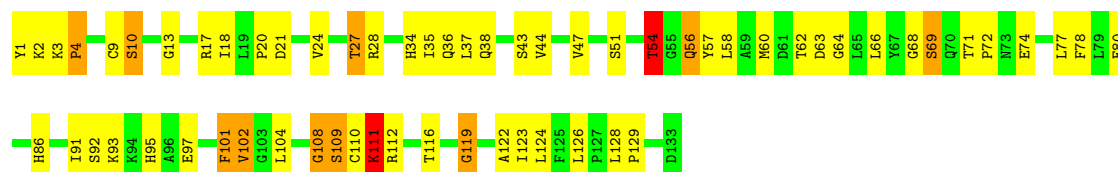
4.2.2 Score per residue for model 2

- Molecule 1: Heparin-binding growth factor 1





- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13

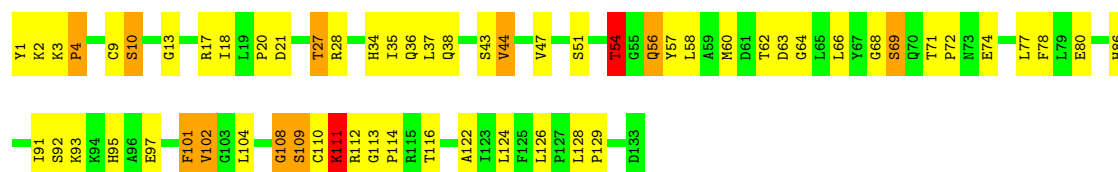


- Molecule 3: Protein S100-A13

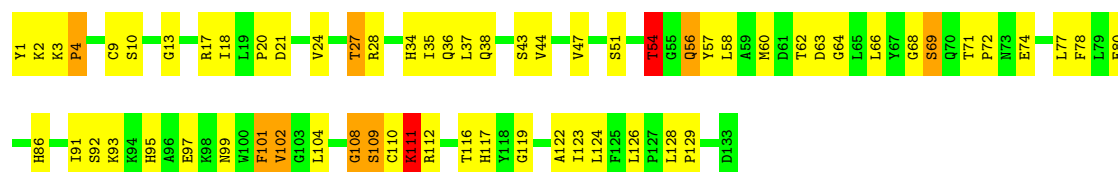


4.2.3 Score per residue for model 3

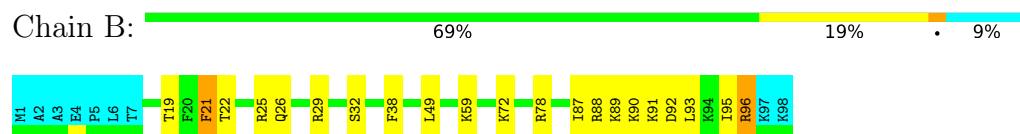
- Molecule 1: Heparin-binding growth factor 1



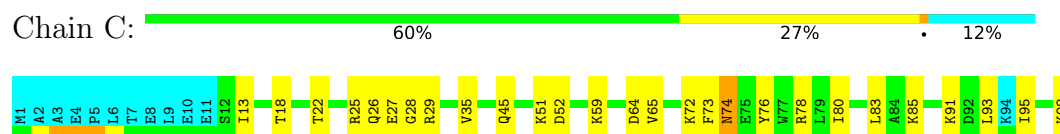
- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13

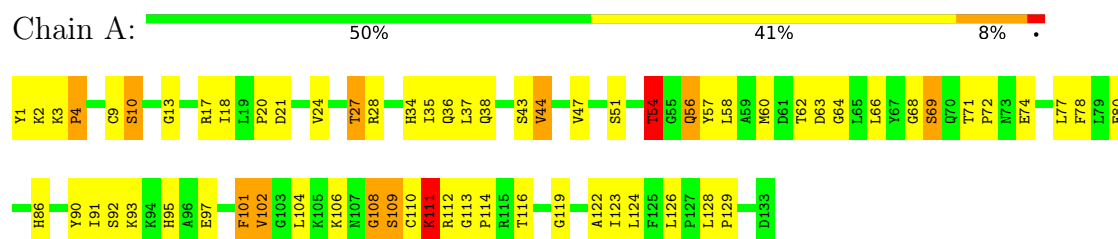


• Molecule 3: Protein S100-A13

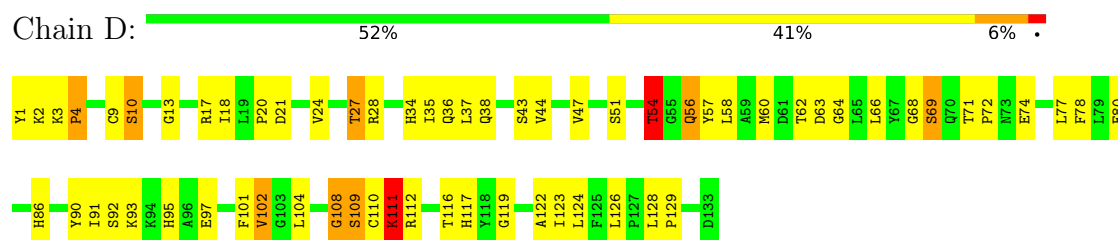


4.2.4 Score per residue for model 4

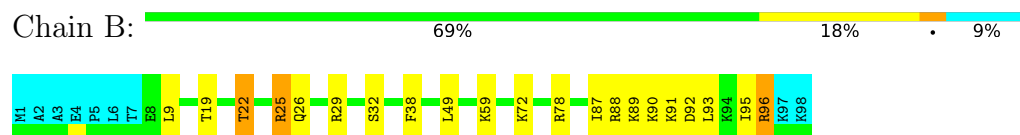
• Molecule 1: Heparin-binding growth factor 1



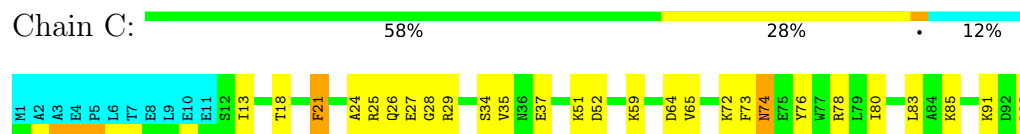
• Molecule 1: Heparin-binding growth factor 1



• Molecule 2: Protein S100-A13

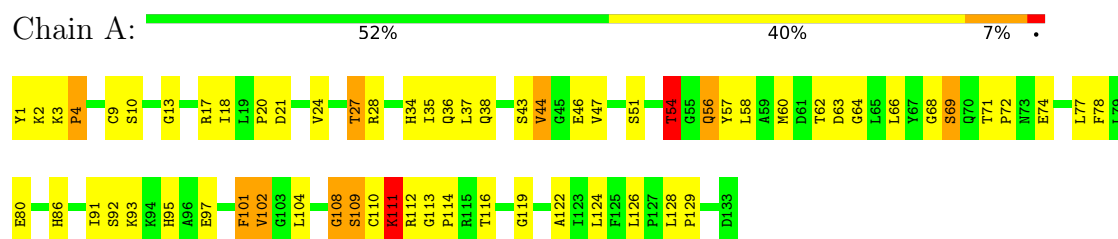


• Molecule 3: Protein S100-A13

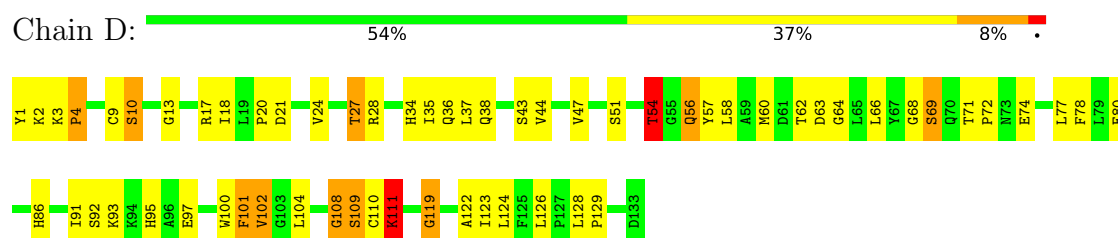


4.2.5 Score per residue for model 5

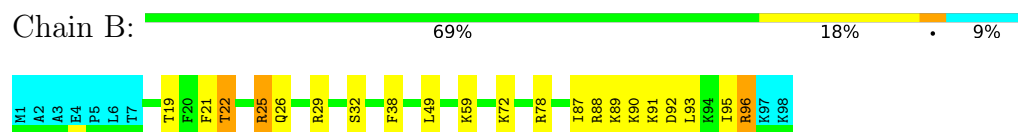
- Molecule 1: Heparin-binding growth factor 1



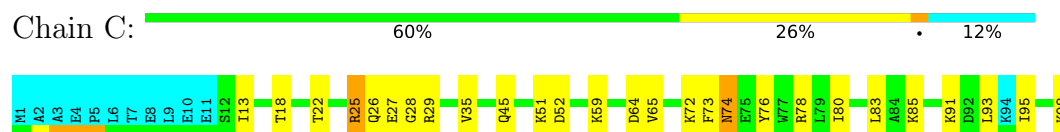
- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13

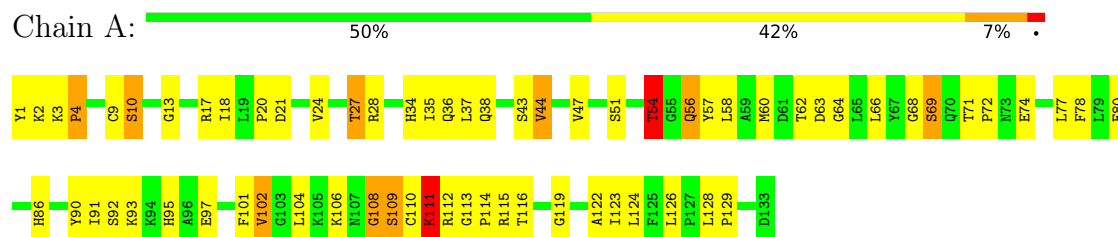


- Molecule 3: Protein S100-A13

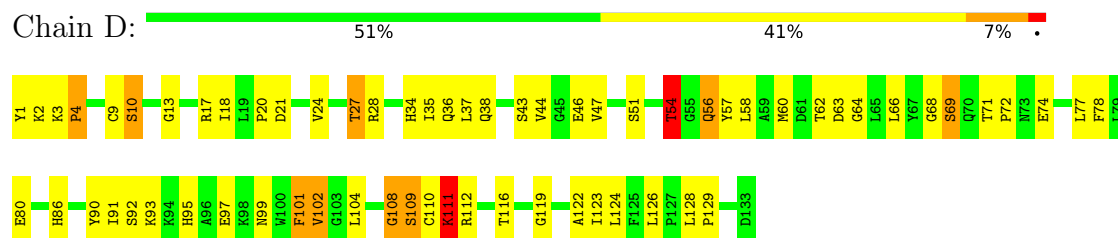


4.2.6 Score per residue for model 6

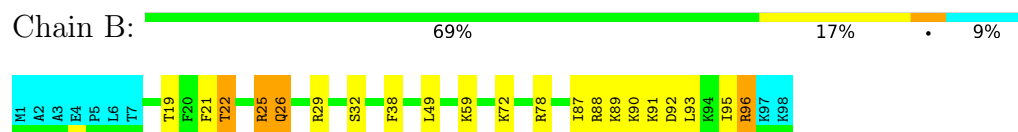
- Molecule 1: Heparin-binding growth factor 1



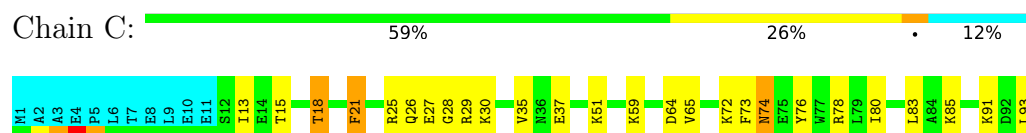
- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13

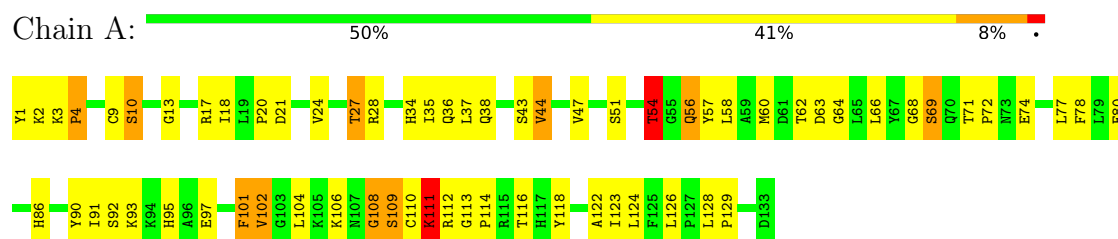


- Molecule 3: Protein S100-A13

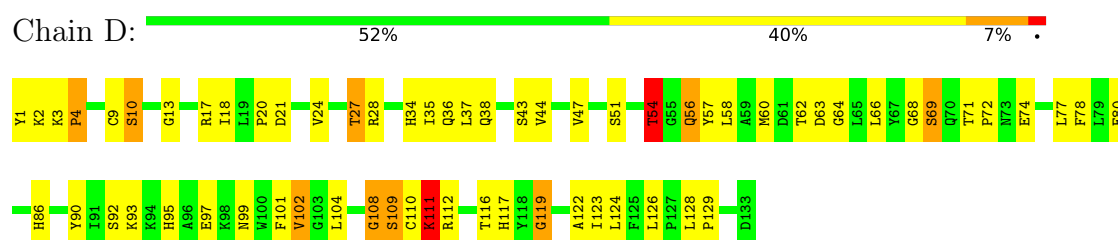


4.2.7 Score per residue for model 7 (medoid)

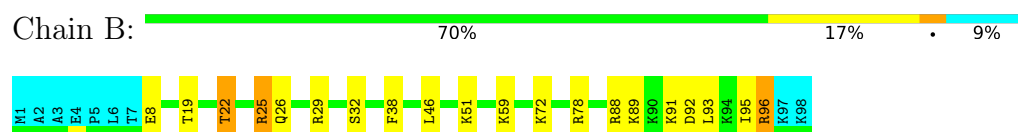
- Molecule 1: Heparin-binding growth factor 1



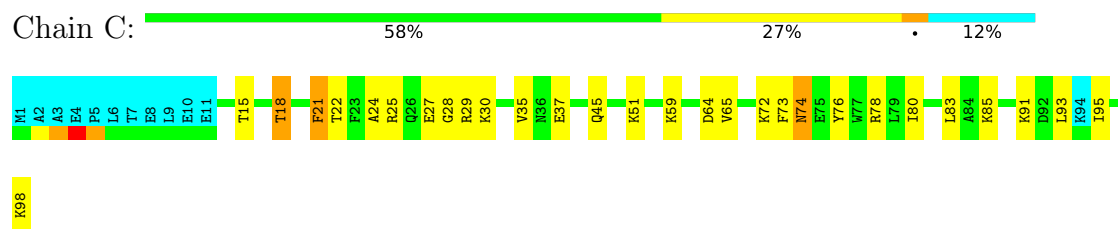
- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13

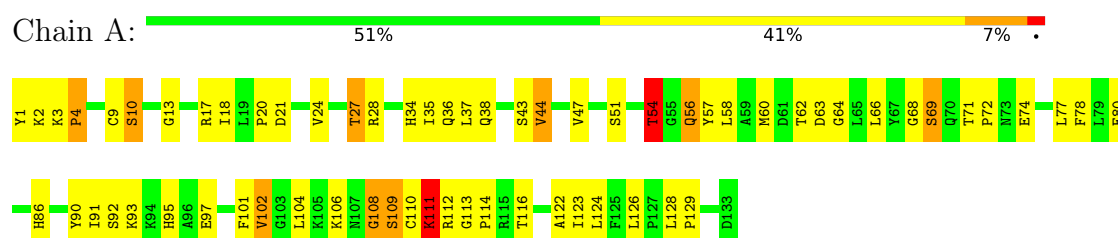


- Molecule 3: Protein S100-A13

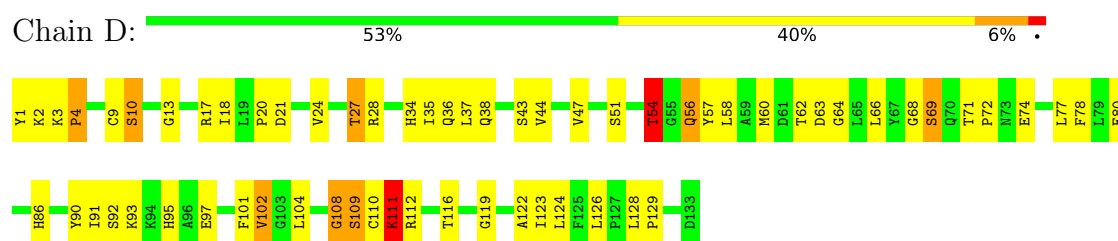


4.2.8 Score per residue for model 8

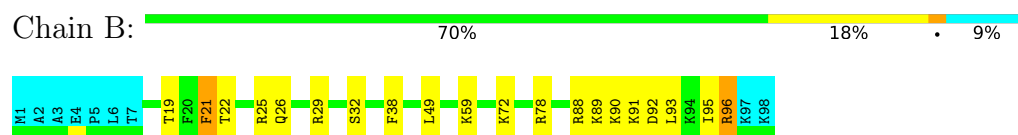
- Molecule 1: Heparin-binding growth factor 1



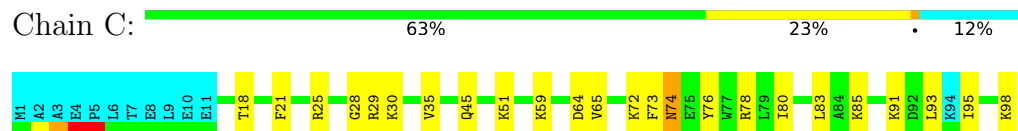
- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13

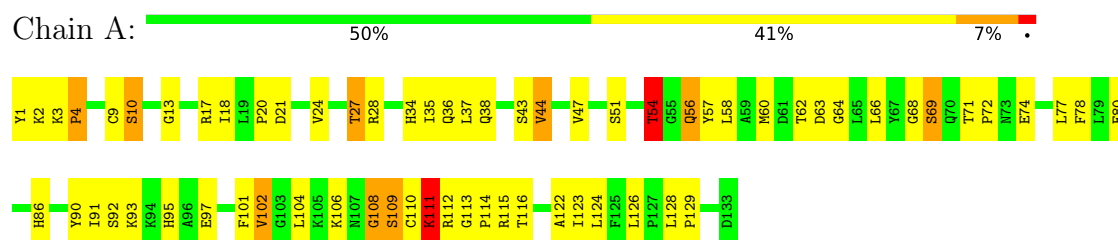


- Molecule 3: Protein S100-A13

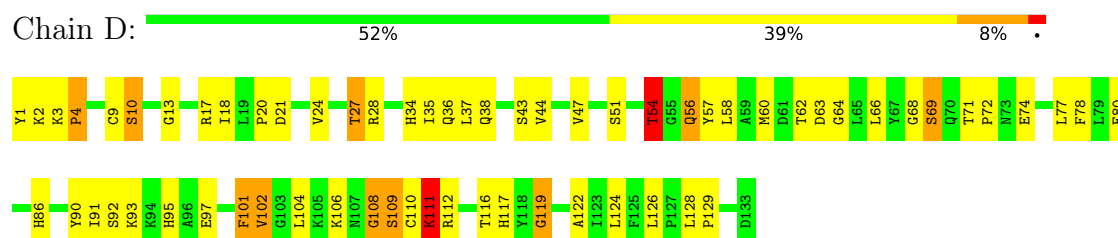


4.2.9 Score per residue for model 9

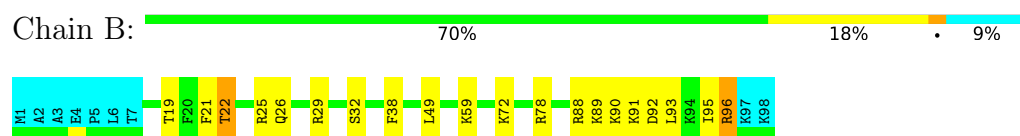
- Molecule 1: Heparin-binding growth factor 1



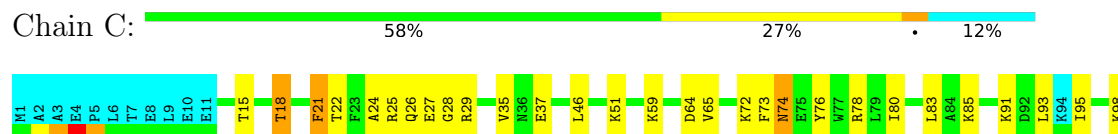
- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13

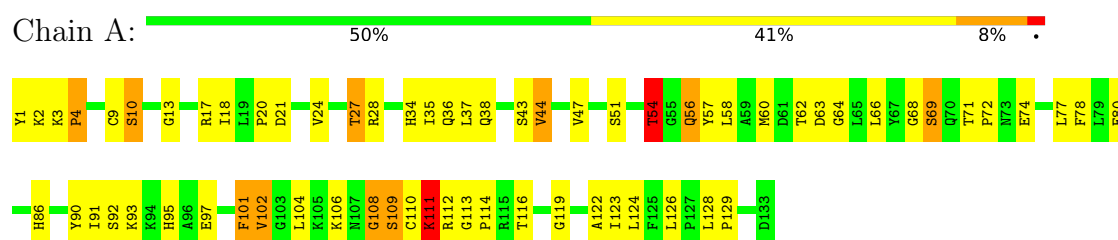


- Molecule 3: Protein S100-A13

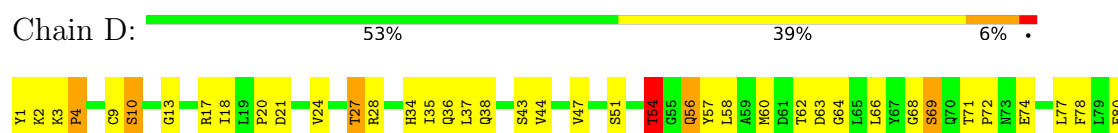


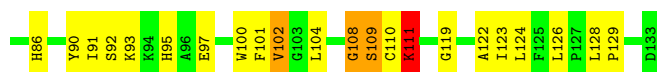
4.2.10 Score per residue for model 10

- Molecule 1: Heparin-binding growth factor 1



- Molecule 1: Heparin-binding growth factor 1





• Molecule 2: Protein S100-A13



• Molecule 3: Protein S100-A13



4.2.11 Score per residue for model 11

• Molecule 1: Heparin-binding growth factor 1



• Molecule 1: Heparin-binding growth factor 1

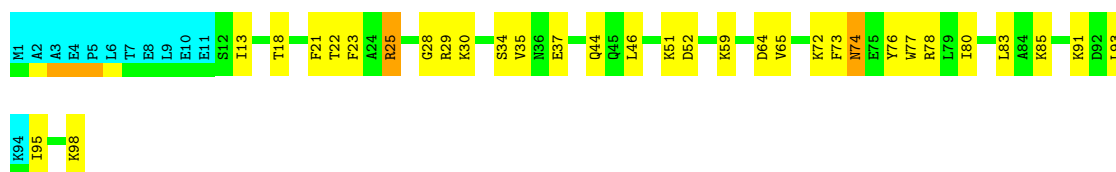


• Molecule 2: Protein S100-A13



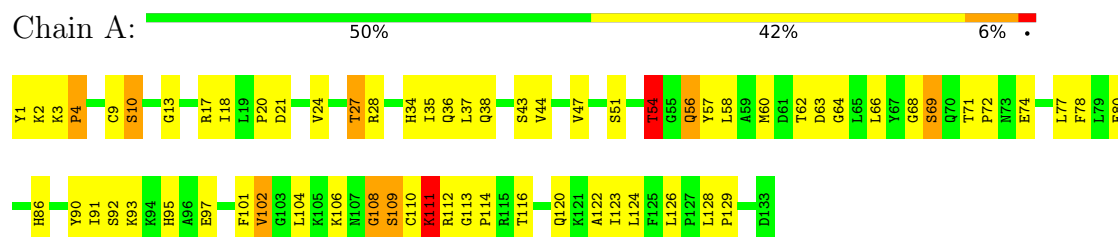
• Molecule 3: Protein S100-A13



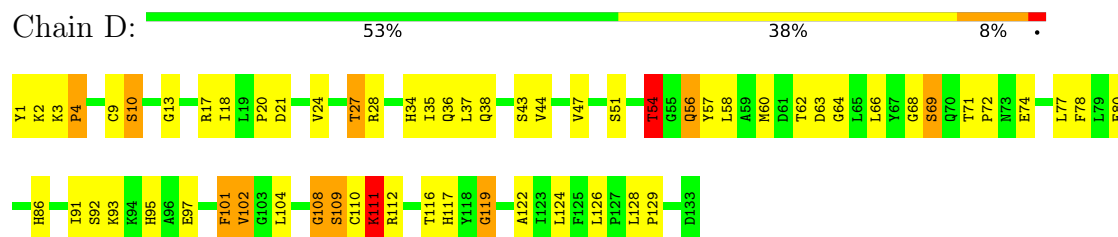


4.2.12 Score per residue for model 12

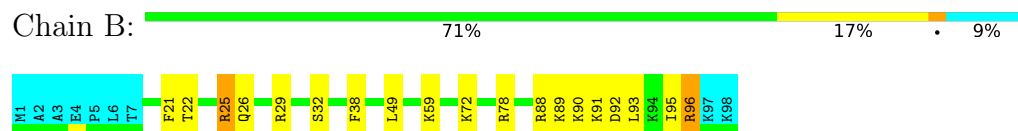
- Molecule 1: Heparin-binding growth factor 1



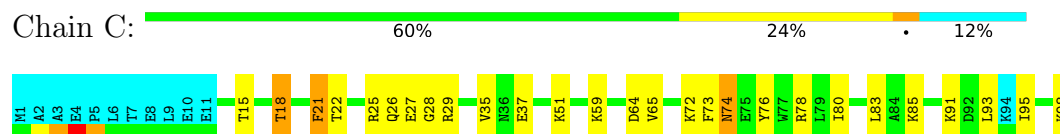
- Molecule 1: Heparin-binding growth factor 1



- Molecule 2: Protein S100-A13



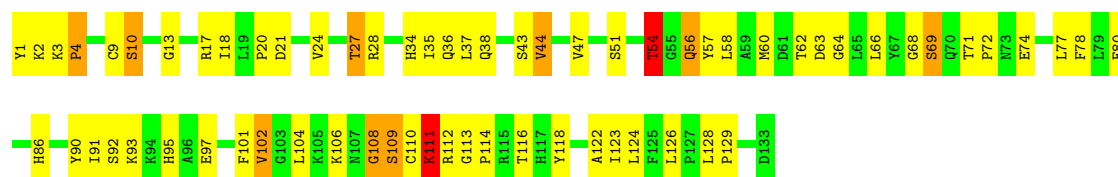
- Molecule 3: Protein S100-A13



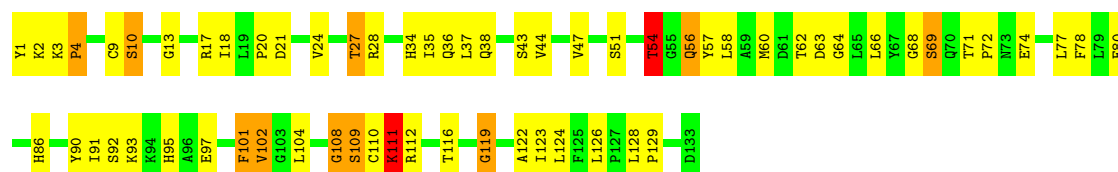
4.2.13 Score per residue for model 13

- Molecule 1: Heparin-binding growth factor 1





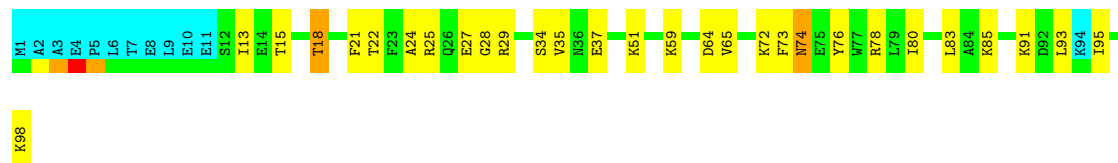
• Molecule 1: Heparin-binding growth factor 1



• Molecule 2: Protein S100-A13

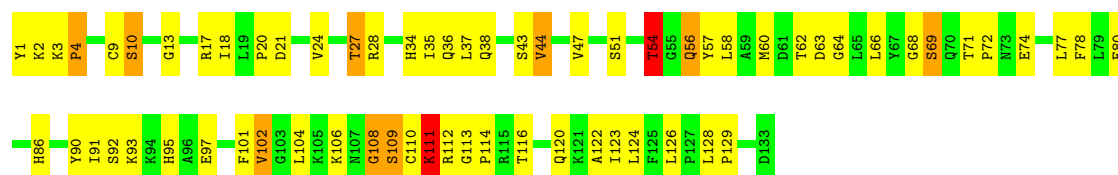


• Molecule 3: Protein S100-A13

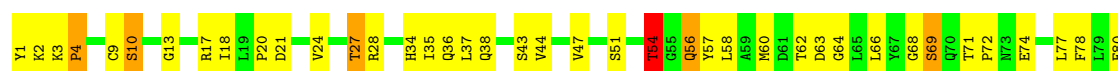


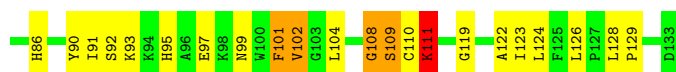
4.2.14 Score per residue for model 14

• Molecule 1: Heparin-binding growth factor 1



• Molecule 1: Heparin-binding growth factor 1





• Molecule 2: Protein S100-A13



• Molecule 3: Protein S100-A13

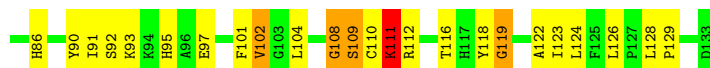


4.2.15 Score per residue for model 15

• Molecule 1: Heparin-binding growth factor 1



• Molecule 1: Heparin-binding growth factor 1

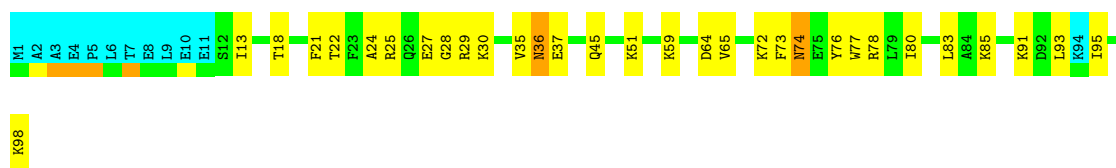


• Molecule 2: Protein S100-A13



• Molecule 3: Protein S100-A13

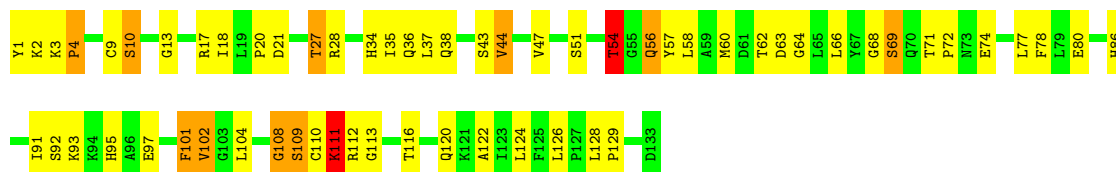




4.2.16 Score per residue for model 16

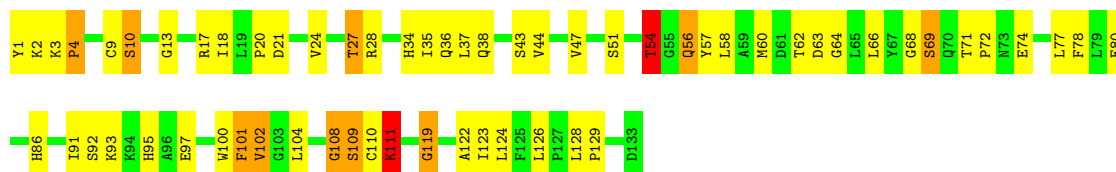
- Molecule 1: Heparin-binding growth factor 1

Chain A: 54% 37% 8% .



- Molecule 1: Heparin-binding growth factor 1

Chain D: 54% 37% 8% .



- Molecule 2: Protein S100-A13

Chain B: 72% 16% 9% .



- Molecule 3: Protein S100-A13

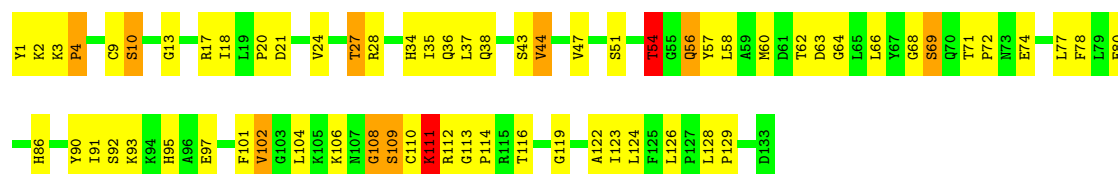
Chain C: 61% 23% 12% .



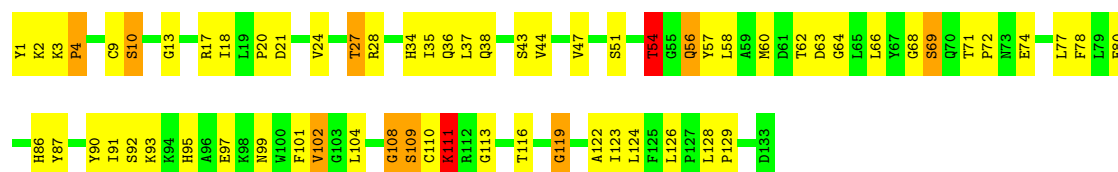
4.2.17 Score per residue for model 17

- Molecule 1: Heparin-binding growth factor 1

Chain A: 50% 41% 7% .



• Molecule 1: Heparin-binding growth factor 1



• Molecule 2: Protein S100-A13

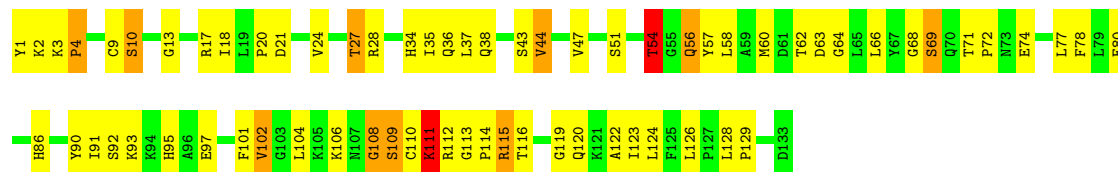


• Molecule 3: Protein S100-A13



4.2.18 Score per residue for model 18

• Molecule 1: Heparin-binding growth factor 1



• Molecule 1: Heparin-binding growth factor 1





• Molecule 2: Protein S100-A13

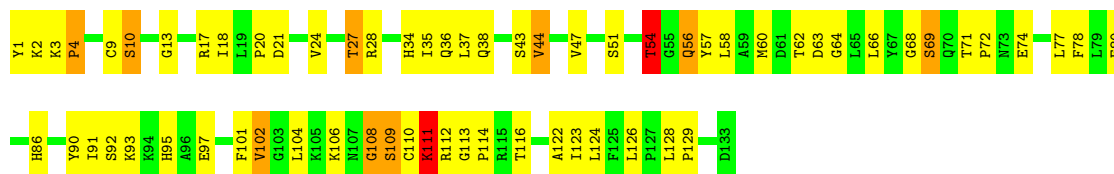


• Molecule 3: Protein S100-A13

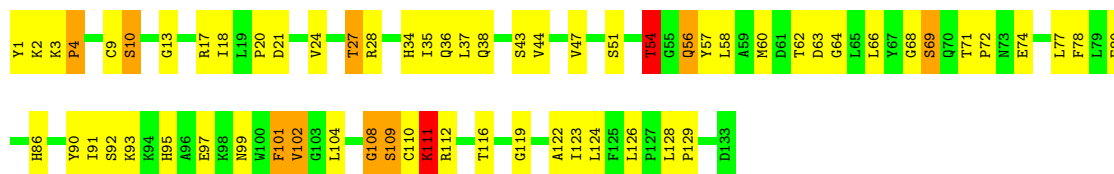


4.2.19 Score per residue for model 19

• Molecule 1: Heparin-binding growth factor 1



• Molecule 1: Heparin-binding growth factor 1



• Molecule 2: Protein S100-A13



• Molecule 3: Protein S100-A13

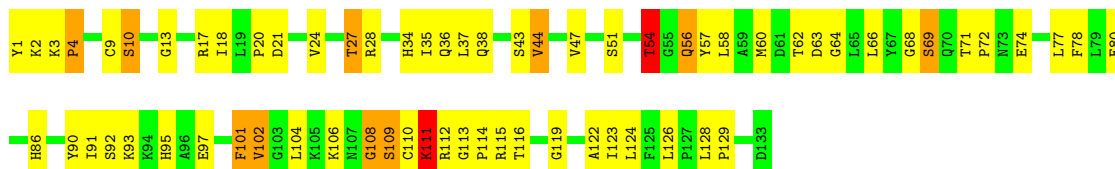




4.2.20 Score per residue for model 20

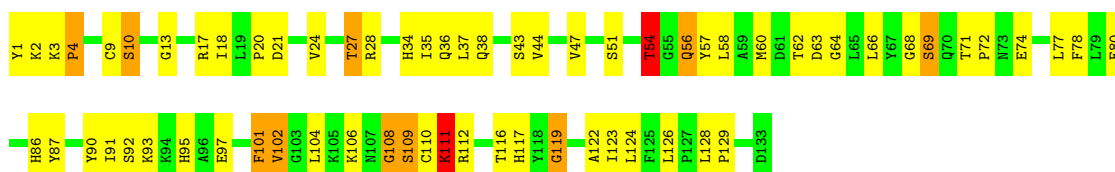
- Molecule 1: Heparin-binding growth factor 1

Chain A: 50% 41% 8% .



- Molecule 1: Heparin-binding growth factor 1

Chain D: 50% 41% 8% .



- Molecule 2: Protein S100-A13

Chain B: 72% 16% 9% .



- Molecule 3: Protein S100-A13

Chain C: 60% 23% 12% .



5 Refinement protocol and experimental data overview ⓘ

The models were refined using the following method: *simulated annealing*.

Of the 5000 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
ARIA	structure solution	1.2 & 2.2
CNS	structure solution	1.2
CNS	refinement	1.2

No chemical shift data was provided.

6 Model quality i

6.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: DLY

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	0.37±0.00	0±0/1088 (0.0± 0.0%)	0.60±0.00	0±0/1466 (0.0± 0.0%)
1	D	0.37±0.00	0±0/1088 (0.0± 0.0%)	0.60±0.00	0±0/1466 (0.0± 0.0%)
2	B	0.58±0.01	0±0/748 (0.0± 0.0%)	1.21±0.01	2±0/1000 (0.2± 0.0%)
3	C	0.58±0.01	0±0/722 (0.0± 0.0%)	1.33±0.01	5±1/961 (0.5± 0.1%)
All	All	0.46	0/72920 (0.0%)	0.93	136/97860 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
2	B	0.0±0.0	4.0±0.0
3	C	0.0±0.0	7.0±0.0
All	All	0	220

There are no bond-length outliers.

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
3	C	80	ILE	CA-C-N	6.40	127.04	120.00	1	20
3	C	80	ILE	C-N-CA	6.40	127.04	120.00	1	20
2	B	95	ILE	CA-C-N	6.38	133.18	121.70	5	20
2	B	95	ILE	C-N-CA	6.38	133.18	121.70	5	20
3	C	73	PHE	CA-C-N	6.36	133.69	121.54	12	20
3	C	73	PHE	C-N-CA	6.36	133.69	121.54	12	20
3	C	26	GLN	N-CA-C	5.79	117.39	111.14	20	2
3	C	27	GLU	CA-C-N	5.12	128.05	121.13	13	7

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
3	C	27	GLU	C-N-CA	5.12	128.05	121.13	13	7

There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
2	B	38	PHE	Sidechain	20
2	B	78	ARG	Sidechain	20
2	B	89	LYS	Peptide	20
2	B	96	ARG	Sidechain	20
3	C	28	GLY	Peptide	20
3	C	35	VAL	Peptide,Mainchain	20
3	C	64	ASP	Peptide	20
3	C	74	ASN	Peptide	20
3	C	76	TYR	Sidechain	20
3	C	78	ARG	Sidechain	20
3	C	91	LYS	Peptide	20

6.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	1064	1049	1046	51±4
1	D	1064	1049	1046	49±2
2	B	739	756	755	4±1
3	C	714	740	739	5±2
All	All	71620	71880	71720	2131

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:66:LEU:HD12	1:A:110:CYS:HB2	0.74	1.59	16	20

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:D:66:LEU:HD12	1:D:110:CYS:HB2	0.69	1.63	15	20
3:C:22:THR:HG21	1:D:119:GLY:HA3	0.68	1.66	20	12
1:A:3:LYS:CB	1:A:4:PRO:HD2	0.67	2.20	9	20
3:C:21:PHE:O	3:C:25:ARG:HG2	0.67	1.89	6	10
1:D:3:LYS:CB	1:D:4:PRO:HD2	0.65	2.20	15	20
2:B:87:ILE:HD13	3:C:13:ILE:HD11	0.63	1.70	4	10
1:A:64:GLY:HA2	1:A:113:GLY:HA3	0.59	1.73	20	20
1:D:9:CYS:SG	1:D:124:LEU:O	0.58	2.58	3	20
2:B:19:THR:O	2:B:22:THR:HG22	0.58	1.98	3	15
1:A:9:CYS:SG	1:A:124:LEU:O	0.58	2.57	3	20
1:D:101:PHE:CD2	1:D:123:ILE:HG21	0.57	2.34	11	17
1:A:90:TYR:O	1:A:102:VAL:HG12	0.56	2.00	7	14
1:A:3:LYS:HB3	1:A:4:PRO:HD2	0.56	1.78	9	20
1:A:60:MET:HG3	1:A:66:LEU:CD2	0.55	2.32	7	20
1:D:3:LYS:HB3	1:D:4:PRO:HD2	0.55	1.78	20	20
1:A:66:LEU:HD13	1:A:102:VAL:HG23	0.55	1.77	15	6
1:A:101:PHE:HD2	1:A:123:ILE:HG21	0.55	1.62	10	14
1:D:108:GLY:O	1:D:109:SER:CB	0.55	2.55	4	20
1:A:108:GLY:O	1:A:109:SER:CB	0.55	2.55	5	20
1:D:60:MET:HG3	1:D:66:LEU:CD2	0.55	2.32	20	20
1:D:58:LEU:HA	1:D:68:GLY:HA3	0.53	1.80	14	20
1:D:20:PRO:HA	1:D:56:GLN:HG3	0.53	1.80	3	20
1:A:58:LEU:HA	1:A:68:GLY:HA3	0.53	1.80	10	20
1:D:60:MET:HG3	1:D:66:LEU:HD22	0.53	1.81	16	20
2:B:49:LEU:HA	2:B:90:LYS:NZ	0.53	2.19	10	16
1:D:101:PHE:HD2	1:D:123:ILE:HG21	0.53	1.62	11	17
1:D:3:LYS:CG	1:D:4:PRO:HD2	0.53	2.34	19	20
1:A:60:MET:HG3	1:A:66:LEU:HD22	0.52	1.81	13	20
2:B:25:ARG:HD2	2:B:25:ARG:C	0.52	2.29	5	20
1:D:34:HIS:HA	1:D:54:THR:HG23	0.52	1.82	16	20
1:D:91:ILE:HA	1:D:101:PHE:HB3	0.52	1.82	10	19
1:A:17:ARG:NH1	1:A:27:THR:OG1	0.52	2.40	8	20
1:A:3:LYS:CG	1:A:4:PRO:HD2	0.52	2.34	20	20
1:A:20:PRO:HA	1:A:56:GLN:HG3	0.52	1.81	4	20
1:A:91:ILE:HA	1:A:101:PHE:HB3	0.52	1.81	12	20
1:D:17:ARG:NH1	1:D:27:THR:OG1	0.52	2.40	15	20
1:A:34:HIS:HA	1:A:54:THR:HG23	0.51	1.82	1	20
1:D:9:CYS:SG	1:D:122:ALA:HA	0.51	2.45	15	20
3:C:24:ALA:HA	3:C:37:GLU:OE1	0.51	2.05	15	6
1:D:86:HIS:CE1	1:D:126:LEU:HD11	0.51	2.41	3	20
1:D:66:LEU:HD13	1:D:102:VAL:HG23	0.51	1.81	2	20

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
3:C:25:ARG:NH1	1:D:100:TRP:HE1	0.51	2.04	16	1
2:B:87:ILE:HD13	3:C:13:ILE:CD1	0.50	2.35	4	9
1:A:10:SER:HB3	1:A:106:LYS:HD3	0.50	1.83	18	14
1:A:112:ARG:O	1:A:116:THR:HG23	0.50	2.07	20	20
1:A:9:CYS:SG	1:A:122:ALA:HA	0.50	2.46	9	20
1:A:86:HIS:CE1	1:A:126:LEU:HD11	0.50	2.42	5	20
3:C:15:THR:O	3:C:18:THR:HG22	0.49	2.07	17	9
3:C:25:ARG:O	3:C:25:ARG:HD2	0.49	2.07	11	2
3:C:25:ARG:HB3	3:C:30:LYS:HB3	0.49	1.84	1	7
3:C:25:ARG:HD3	1:D:117:HIS:CG	0.49	2.42	20	1
1:A:10:SER:HB3	1:A:106:LYS:HB2	0.48	1.85	7	14
1:D:128:LEU:N	1:D:129:PRO:CD	0.48	2.76	4	20
1:A:128:LEU:N	1:A:129:PRO:CD	0.48	2.76	2	20
1:D:101:PHE:CZ	1:D:117:HIS:HA	0.48	2.44	4	4
1:A:13:GLY:O	1:A:28:ARG:HB3	0.48	2.09	5	20
1:D:101:PHE:CE2	1:D:123:ILE:HD13	0.48	2.43	20	8
1:D:24:VAL:HG13	1:D:110:CYS:SG	0.48	2.47	5	20
1:D:128:LEU:N	1:D:129:PRO:HD3	0.48	2.24	17	20
1:A:128:LEU:N	1:A:129:PRO:HD3	0.48	2.24	2	20
1:D:13:GLY:O	1:D:28:ARG:HB3	0.48	2.09	2	20
1:D:56:GLN:OE1	1:D:56:GLN:N	0.48	2.47	4	20
1:A:56:GLN:N	1:A:56:GLN:OE1	0.48	2.47	2	20
1:D:104:LEU:C	1:D:104:LEU:HD12	0.48	2.34	15	20
1:A:104:LEU:C	1:A:104:LEU:HD12	0.47	2.34	5	20
3:C:22:THR:HG21	1:D:119:GLY:N	0.47	2.24	14	1
1:A:3:LYS:CB	1:A:4:PRO:CD	0.47	2.93	9	20
1:D:63:ASP:HB2	1:D:95:HIS:CD2	0.47	2.45	3	20
1:A:63:ASP:HB2	1:A:95:HIS:CD2	0.47	2.45	1	20
1:A:115:ARG:HG3	2:B:26:GLN:OE1	0.47	2.09	20	3
1:A:77:LEU:CD1	1:A:93:LYS:HB3	0.47	2.40	13	20
1:D:1:TYR:CE2	1:D:3:LYS:HD2	0.47	2.45	12	20
1:D:77:LEU:CD1	1:D:93:LYS:HB3	0.47	2.40	20	20
1:A:108:GLY:O	1:A:109:SER:HB2	0.46	2.11	14	20
1:D:47:VAL:CG1	1:D:80:GLU:HB2	0.46	2.41	16	20
1:A:1:TYR:CE2	1:A:3:LYS:HD2	0.46	2.45	5	20
1:A:64:GLY:CA	1:A:113:GLY:HA3	0.46	2.40	20	12
1:D:90:TYR:O	1:D:102:VAL:HG12	0.46	2.10	19	14
1:D:112:ARG:O	1:D:116:THR:HG23	0.46	2.10	7	15
1:A:66:LEU:HD13	1:A:102:VAL:CG2	0.46	2.40	15	5
1:A:3:LYS:O	1:A:4:PRO:C	0.46	2.59	2	20
1:D:3:LYS:O	1:D:4:PRO:C	0.46	2.59	5	20

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:57:TYR:HB3	1:A:72:PRO:HB3	0.46	1.87	2	20
1:D:57:TYR:HB3	1:D:72:PRO:HB3	0.46	1.88	8	20
1:D:108:GLY:O	1:D:109:SER:HB2	0.46	2.11	15	20
2:B:21:PHE:O	2:B:25:ARG:HG3	0.46	2.11	18	5
1:A:66:LEU:CD1	1:A:110:CYS:HB2	0.46	2.40	17	11
1:A:47:VAL:CG1	1:A:80:GLU:HB2	0.46	2.41	11	20
1:D:9:CYS:SG	1:D:122:ALA:O	0.46	2.74	7	20
1:D:28:ARG:HG3	1:D:28:ARG:O	0.45	2.12	11	20
1:D:3:LYS:HB3	1:D:4:PRO:CD	0.45	2.41	4	20
1:D:110:CYS:C	1:D:111:LYS:HD2	0.45	2.37	6	20
1:A:60:MET:HE2	1:A:78:PHE:CZ	0.45	2.47	6	20
1:A:110:CYS:C	1:A:111:LYS:HD2	0.45	2.37	2	20
1:D:109:SER:HB3	1:D:111:LYS:HD3	0.45	1.88	15	20
1:A:64:GLY:HA2	1:A:114:PRO:HD3	0.45	1.88	7	18
1:D:60:MET:HE2	1:D:78:PHE:CZ	0.45	2.47	7	20
1:A:101:PHE:CD2	1:A:123:ILE:HG21	0.45	2.47	9	14
1:A:120:GLN:HG2	2:B:25:ARG:NH2	0.45	2.27	16	1
1:A:3:LYS:HB3	1:A:4:PRO:CD	0.45	2.41	1	20
1:A:109:SER:HB3	1:A:111:LYS:HD3	0.45	1.88	17	20
3:C:22:THR:HG21	1:D:119:GLY:CA	0.45	2.42	3	3
1:D:3:LYS:CB	1:D:4:PRO:CD	0.45	2.93	6	16
1:A:28:ARG:O	1:A:28:ARG:HG3	0.44	2.12	10	20
3:C:21:PHE:CZ	3:C:30:LYS:HD2	0.44	2.48	11	1
1:D:24:VAL:CG1	1:D:110:CYS:SG	0.44	3.06	5	17
2:B:87:ILE:CD1	3:C:13:ILE:HD11	0.44	2.40	4	1
1:A:4:PRO:HB3	1:A:36:GLN:CB	0.44	2.43	6	20
1:A:63:ASP:O	1:A:64:GLY:C	0.44	2.61	1	20
1:A:101:PHE:HD1	1:A:123:ILE:HG21	0.44	1.72	8	1
1:A:9:CYS:SG	1:A:122:ALA:O	0.44	2.76	11	16
1:D:63:ASP:O	1:D:64:GLY:C	0.44	2.61	8	20
1:A:24:VAL:HG13	1:A:110:CYS:SG	0.43	2.53	10	17
1:A:120:GLN:H	2:B:25:ARG:HH12	0.43	1.54	16	1
1:D:68:GLY:O	1:D:69:SER:HB2	0.43	2.13	2	20
1:A:120:GLN:N	2:B:25:ARG:HH12	0.43	2.11	16	1
1:D:4:PRO:HB3	1:D:36:GLN:CB	0.43	2.43	6	20
3:C:34:SER:O	3:C:37:GLU:HB2	0.43	2.12	10	5
3:C:25:ARG:NH2	1:D:100:TRP:HE1	0.43	2.11	5	1
3:C:26:GLN:HB3	3:C:37:GLU:HG2	0.43	1.90	6	4
1:A:10:SER:CB	1:A:106:LYS:HD3	0.43	2.44	7	4
1:A:68:GLY:O	1:A:69:SER:HB2	0.43	2.13	7	20
1:A:101:PHE:O	1:A:102:VAL:C	0.43	2.61	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:24:VAL:CG1	1:A:110:CYS:SG	0.43	3.06	9	14
2:B:81:GLY:HA2	3:C:77:TRP:CH2	0.42	2.50	15	2
3:C:25:ARG:NH2	3:C:26:GLN:HA	0.42	2.29	17	1
3:C:25:ARG:HD2	3:C:25:ARG:C	0.42	2.40	3	1
1:A:46:GLU:OE2	1:A:93:LYS:NZ	0.42	2.52	5	1
1:D:113:GLY:HA2	1:D:116:THR:HG23	0.42	1.89	17	1
1:D:20:PRO:HA	1:D:56:GLN:CG	0.42	2.45	3	20
1:A:20:PRO:HA	1:A:56:GLN:CG	0.42	2.45	15	20
1:D:56:GLN:HB2	1:D:69:SER:HA	0.42	1.92	3	20
1:A:101:PHE:CD1	1:A:123:ILE:HG21	0.42	2.50	8	1
1:A:56:GLN:HB2	1:A:69:SER:HA	0.42	1.92	3	20
1:D:46:GLU:OE2	1:D:93:LYS:NZ	0.42	2.52	6	1
1:D:66:LEU:CD1	1:D:110:CYS:HB2	0.41	2.42	15	2
1:A:91:ILE:HA	1:A:101:PHE:CB	0.41	2.45	14	6
3:C:25:ARG:HH12	1:D:100:TRP:CD1	0.41	2.33	10	1
1:A:120:GLN:HG2	2:B:25:ARG:HH22	0.41	1.74	12	3
1:D:77:LEU:HD11	1:D:93:LYS:HB3	0.41	1.92	13	20
2:B:9:LEU:HD13	3:C:46:LEU:HD11	0.41	1.92	1	1
1:A:104:LEU:O	1:A:122:ALA:HB1	0.41	2.15	18	3
1:D:9:CYS:SG	1:D:10:SER:N	0.41	2.94	2	19
3:C:21:PHE:O	3:C:25:ARG:HG3	0.41	2.16	7	1
1:D:10:SER:HB3	1:D:106:LYS:HB2	0.41	1.92	20	2
1:D:87:TYR:CB	1:D:124:LEU:HB3	0.40	2.46	20	3
1:A:9:CYS:SG	1:A:10:SER:N	0.40	2.94	7	14
1:A:44:VAL:O	1:A:44:VAL:HG13	0.40	2.16	15	18
1:A:118:TYR:O	2:B:25:ARG:HG2	0.40	2.16	13	2
3:C:23:PHE:O	3:C:37:GLU:HG2	0.40	2.17	11	1

6.3 Torsion angles ⓘ

6.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the backbone conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	131/133 (98%)	95±1 (73±0%)	26±1 (20±1%)	10±0 (7±0%)	1	15
1	D	131/133 (98%)	96±0 (73±0%)	25±1 (19±0%)	10±0 (8±0%)	1	14

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	89/98 (91%)	68±1 (76±1%)	17±1 (19±1%)	4±0 (5±0%)	3	25
3	C	85/98 (87%)	65±1 (76±1%)	16±1 (18±1%)	5±1 (5±1%)	2	21
All	All	8720/9240 (94%)	6482 (74%)	1673 (19%)	565 (6%)	2	18

All 32 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	4	PRO	20
1	A	10	SER	20
1	A	44	VAL	20
1	A	54	THR	20
1	A	69	SER	20
1	A	102	VAL	20
1	A	108	GLY	20
1	A	109	SER	20
1	A	111	LYS	20
2	B	29	ARG	20
2	B	92	ASP	20
2	B	93	LEU	20
2	B	96	ARG	20
3	C	29	ARG	20
3	C	65	VAL	20
3	C	74	ASN	20
3	C	95	ILE	20
1	D	4	PRO	20
1	D	10	SER	20
1	D	44	VAL	20
1	D	54	THR	20
1	D	69	SER	20
1	D	102	VAL	20
1	D	108	GLY	20
1	D	109	SER	20
1	D	111	LYS	20
1	D	119	GLY	18
1	A	119	GLY	10
3	C	27	GLU	9
2	B	51	LYS	4
3	C	26	GLN	3
3	C	36	ASN	1

6.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the sidechain conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	117/117 (100%)	99±0 (85±0%)	18±0 (15±0%)	5	42
1	D	117/117 (100%)	99±1 (84±1%)	18±1 (16±1%)	5	41
2	B	84/89 (94%)	76±1 (90±1%)	8±1 (10±1%)	10	54
3	C	81/90 (90%)	71±1 (88±1%)	10±1 (12±1%)	7	48
All	All	7980/8260 (97%)	6900 (86%)	1080 (14%)	6	45

All 68 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	2	LYS	20
1	A	18	ILE	20
1	A	21	ASP	20
1	A	27	THR	20
1	A	35	ILE	20
1	A	37	LEU	20
1	A	38	GLN	20
1	A	43	SER	20
1	A	51	SER	20
1	A	54	THR	20
1	A	56	GLN	20
1	A	62	THR	20
1	A	71	THR	20
1	A	74	GLU	20
1	A	92	SER	20
1	A	97	GLU	20
1	A	111	LYS	20
2	B	26	GLN	20
2	B	32	SER	20
2	B	59	LYS	20
2	B	72	LYS	20
2	B	88	ARG	20
2	B	91	LYS	20
3	C	18	THR	20
3	C	59	LYS	20

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Mol	Chain	Res	Type	Models (Total)
3	C	72	LYS	20
3	C	83	LEU	20
3	C	85	LYS	20
3	C	93	LEU	20
3	C	98	LYS	20
1	D	2	LYS	20
1	D	18	ILE	20
1	D	21	ASP	20
1	D	27	THR	20
1	D	35	ILE	20
1	D	37	LEU	20
1	D	38	GLN	20
1	D	43	SER	20
1	D	51	SER	20
1	D	54	THR	20
1	D	56	GLN	20
1	D	62	THR	20
1	D	71	THR	20
1	D	74	GLU	20
1	D	92	SER	20
1	D	97	GLU	20
1	D	111	LYS	20
3	C	51	LYS	19
2	B	22	THR	14
2	B	25	ARG	13
1	D	101	PHE	13
2	B	21	PHE	13
1	A	101	PHE	11
3	C	21	PHE	11
1	D	99	ASN	7
3	C	45	GLN	6
3	C	52	ASP	6
3	C	25	ARG	6
3	C	26	GLN	4
2	B	46	LEU	3
2	B	8	GLU	3
3	C	46	LEU	3
1	D	117	HIS	2
1	A	115	ARG	2
2	B	9	LEU	1
3	C	44	GLN	1
3	C	36	ASN	1

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Mol	Chain	Res	Type	Models (Total)
1	D	118	TYR	1

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

3 non-standard protein/DNA/RNA residues are modelled in this entry.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided